

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE110118\
 Data File : BE098143.D
 Acq On : 1 Nov 2018 13:04
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE110118

Quant Time: Nov 01 13:34:12 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 13:00:52 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	103	0.00
2	1,4-Dioxane	0.400	0.410	-2.5	106	0.00
3	n-Nitrosodimethylamine	0.400	0.423	-5.7	117	0.00
4 S	2-Fluorophenol	0.400	0.379	5.3	95	0.00
5 S	Phenol-d6	0.400	0.384	4.0	99	0.00
6	bis(2-Chloroethyl)ether	0.400	0.379	5.3	101	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	102	0.01
8 S	Nitrobenzene-d5	0.400	0.391	2.3	103	0.00
9	Naphthalene	0.400	0.387	3.3	101	0.00
10	Hexachlorobutadiene	0.400	0.399	0.3	105	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.386	3.5	99	0.00
12	2-Methylnaphthalene	0.400	0.382	4.5	101	0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	97	0.00
14 S	2,4,6-Tribromophenol	0.400	0.339	15.3	98	0.00
15 S	2-Fluorobiphenyl	0.400	0.373	6.8	100	0.00
16	Acenaphthylene	0.400	0.378	5.5	99	0.00
17	Acenaphthene	0.400	0.394	1.5	100	0.00
18	Fluorene	0.400	0.380	5.0	99	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	102	0.00
20	4-Bromophenyl-phenylether	0.400	0.374	6.5	98	0.00
21	Hexachlorobenzene	0.400	0.380	5.0	101	0.00
22	Pentachlorophenol	0.400	0.349	12.8	82	0.00
23	Phenanthrene	0.400	0.372	7.0	99	0.00
24	Anthracene	0.400	0.375	6.3	101	0.00
25 SURR	Fluoranthene-d10	0.400	0.365	8.8	98	0.00
26	Fluoranthene	0.400	0.369	7.8	100	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	99	0.00
28	Pvrene	0.400	0.372	7.0	100	0.00
29 S	Terphenyl-d14	0.400	0.362	9.5	98	0.00
30	Benzo(a)anthracene	0.400	0.368	8.0	99	0.00
31	Chrysene	0.400	0.370	7.5	99	0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.390	2.5	104	0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.375	6.3	100	0.00
34 I	Perylene-d12	0.400	0.400	0.0	99	0.00
35	Benzo(b)fluoranthene	0.400	0.377	5.8	100	0.00
36	Benzo(k)fluoranthene	0.400	0.377	5.8	101	0.00
37 C	Benzo(a)pyrene	0.400	0.371	7.3	98	0.00
38	Dibenzo(a,h)anthracene	0.400	0.371	7.3	97	0.00
39	Benzo(a,h,i)perylene	0.400	0.370	7.5	97	0.00

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Compound	Amount	Calc.	%Dev Area	% Dev(min)

(#) = Out of Range	SPCC's out = 0 CCC's out = 0			